

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
 Data File : BG051836.D
 Acq On : 28 Dec 2021 19:29
 Operator : CG/JU
 Sample : PB141709BL
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK709

Quant Time: Dec 29 00:19:57 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 23 13:37:07 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.270	152	34768	20.000	ng/u1	-0.02
20) Naphthalene-d8	11.113	136	146619	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.908	164	101018	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.651	188	224496	20.000	ng/u1	-0.01
79) Chrysene-d12	21.975	240	191117	20.000	ng/u1	#-0.01
88) Perylene-d12	25.464	264	132107	20.000	ng/u1	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.599	96	4868	5.815	ng/uL	0.00
4) Pyridine-d5	4.028	84	80037	31.642	ng/u1	-0.02
7) Phenol-d5	7.412	99	104588	33.691	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.576	67	62732	33.993	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.800	132	72136	33.411	ng/u1	-0.01
15) 4-Methylphenol-d8	8.969	113	81224	33.849	ng/u1	-0.01
21) Nitrobenzene-d5	9.444	128	37341	34.737	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.173	143	43185	36.750	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.719	165	81713	32.066	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.230	131	98617	29.393	ng/u1	-0.01
46) Dimethylphthalate-d6	14.297	166	258082	31.919	ng/u1	-0.02
49) Acenaphthylene-d8	14.602	160	321096	31.976	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.078	143	25517	19.407	ng/u1	0.00
60) Fluorene-d10	15.895	176	219832	30.347	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	16.000	200	25739	22.006	ng/u1	-0.01
73) Anthracene-d10	17.751	188	355033	33.152	ng/u1	-0.01
81) Pyrene-d10	20.030	212	405618	35.264	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.223	264	236403	34.010	ng/u1	-0.03

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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